

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102517\
 Data File : BF099809.D
 Acq On : 25 Oct 2017 13:04
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 25 17:02:42 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 25 17:02:52 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	82273	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	343755	20.00	ng	0.00
38) Acenaphthene-d10	9.84	164	161134	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	295943	20.00	ng	0.00
75) Chrysene-d12	13.98	240	236626	20.00	ng	0.00
86) Perylene-d12	15.44	264	219902	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.41	112	432603	82.55	ng	0.00
7) Phenol-d6	6.44	99	503001	80.95	ng	0.00
23) Nitrobenzene-d5	7.37	82	413925	77.52	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	116902	79.61	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	851985	81.58	ng	0.00
78) Terphenyl-d14	12.92	244	812409	75.03	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.50	88	92732	40.072	ng	# 89
3) Pyridine	3.28	79	218407	39.246	ng	89
4) n-Nitrosodimethylamine	3.23	42	82366	41.429	ng	# 71
6) Aniline	6.46	93	328617	40.788	ng	98
8) 2-Chlorophenol	6.58	128	226015	40.467	ng	95
9) Benzaldehyde	6.35	77	148439	39.977	ng	93
10) Phenol	6.45	94	276076	40.467	ng	97
11) bis(2-Chloroethyl)ether	6.53	93	215188	40.846	ng	95
12) 1,3-Dichlorobenzene	6.73	146	250869	40.004	ng	96
13) 1,4-Dichlorobenzene	6.81	146	258573	40.626	ng	98
14) 1,2-Dichlorobenzene	6.96	146	239339	41.260	ng	97
15) Benzyl Alcohol	6.95	79	160299	40.565	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.07	45	234971	40.151	ng	69
17) 2-Methylphenol	7.06	107	190420	40.759	ng	99
18) Hexachloroethane	7.30	117	83870	39.498	ng	99
19) n-Nitroso-di-n-propylamine	7.21	70	145484	39.953	ng	91
20) 3+4-Methylphenols	7.21	107	224111	41.198	ng	# 61
22) Acetophenone	7.21	105	305899	39.083	ng	# 98
24) Nitrobenzene	7.39	77	210101	39.053	ng	91
25) Isophorone	7.62	82	376019	38.810	ng	96
26) 2-Nitrophenol	7.70	139	125360	40.683	ng	87
27) 2,4-Dimethylphenol	7.74	122	180241	39.699	ng	93
28) bis(2-Chloroethoxy)methane	7.83	93	267574	39.433	ng	99
29) 2,4-Dichlorophenol	7.95	162	189621	40.205	ng	95
30) 1,2,4-Trichlorobenzene	8.02	180	197879	39.267	ng	98
31) Naphthalene	8.10	128	660293	39.793	ng	99
32) Benzoic acid	7.89	122	162903	40.764	ng	97
33) 4-Chloroaniline	8.16	127	281004	41.011	ng	# 94
34) Hexachlorobutadiene	8.21	225	101432	39.132	ng	98
35) Caprolactam	8.55	113	60053	40.489	ng	# 75
36) 4-Chloro-3-methylphenol	8.65	107	188616	39.181	ng	92
37) 2-Methylnaphthalene	8.79	142	445358	40.050	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.96	216	204562	39.307	ng	# 97
40) Hexachlorocyclopentadiene	8.94	237	29834	38.471	ng	96

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102517\
 Data File : BF099809.D
 Acq On : 25 Oct 2017 13:04
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Oct 25 17:02:42 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 25 17:02:52 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.08	196	125428	39.844	nq	99
43) 2,4,5-Trichlorophenol	9.13	196	134401	40.233	nq	# 93
45) 1,1'-Biphenyl	9.26	154	570803	39.158	nq	98
46) 2-Chloronaphthalene	9.29	162	412996	39.012	nq	96
47) 2-Nitroaniline	9.39	65	109492	39.408	nq	84
48) Acenaphthylene	9.70	152	644757	39.544	nq	100
49) Dimethylphthalate	9.56	163	489969	38.397	nq	99
50) 2,6-Dinitrotoluene	9.63	165	105179	39.209	nq	# 82
51) Acenaphthene	9.87	154	388421	38.987	nq	98
52) 3-Nitroaniline	9.81	138	128090	40.524	nq	83
53) 2,4-Dinitrophenol	9.93	184	35696	37.037	nq	# 81
54) Dibenzofuran	10.05	168	545118	38.762	nq	98
55) 4-Nitrophenol	9.98	139	66541	40.290	nq	82
56) 2,4-Dinitrotoluene	10.05	165	126614	39.193	nq	# 94
57) Fluorene	10.39	166	458042	39.338	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.17	232	94604	40.391	nq	93
59) Diethylphthalate	10.26	149	484095	38.848	nq	98
60) 4-Chlorophenyl-phenylether	10.37	204	215334	39.295	nq	95
61) 4-Nitroaniline	10.43	138	124654	41.336	nq	85
62) Azobenzene	10.54	77	408531	39.109	nq	89
64) 4,6-Dinitro-2-methylphenol	10.46	198	73816	39.679	nq	83
65) n-Nitrosodiphenylamine	10.50	169	430175	39.377	nq	99
66) 4-Bromophenyl-phenylether	10.87	248	122469	39.309	nq	# 86
67) Hexachlorobenzene	10.94	284	122551	38.449	nq	# 92
68) Atrazine	11.03	200	126948	38.685	nq	98
69) Pentachlorophenol	11.14	266	53749	39.464	nq	96
70) Phenanthrene	11.36	178	652308	39.057	nq	98
71) Anthracene	11.40	178	662976	39.107	nq	99
72) Carbazole	11.57	167	634626	39.808	nq	99
73) Di-n-butylphthalate	11.88	149	777062	38.215	nq	99
74) Fluoranthene	12.55	202	690631	39.791	nq	99
76) Benzidine	12.67	184	353762	34.166	nq	98
77) Pyrene	12.78	202	698059	38.796	nq	99
79) Butylbenzylphthalate	13.39	149	334235	37.723	nq	87
80) Benzo(a)anthracene	13.97	228	587632	39.174	nq	99
81) 3,3'-Dichlorobenzidine	13.93	252	217411	39.928	nq	99
82) Chrysene	14.01	228	546116	38.725	nq	98
83) Bis(2-ethylhexyl)phthalate	13.94	149	450253	36.186	nq	98
84) Di-n-octyl phthalate	14.55	149	787894	36.894	nq	# 94
85) Indeno(1,2,3-cd)pyrene	16.93	276	483745	37.365	nq	97
87) Benzo(b)fluoranthene	15.02	252	521025	37.522	nq	97
88) Benzo(k)fluoranthene	15.05	252	541767	41.376	nq	99
89) Benzo(a)pyrene	15.39	252	494740	39.664	nq	# 97
90) Dibenzo(a,h)anthracene	16.93	278	409926	36.986	nq	97
91) Benzo(g,h,i)perylene	17.37	276	396955	36.608	nq	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

